

Pietro Sormanni

List of Publications by Citations

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

56

papers

1,838

citations

24

h-index

42

g-index

61

ext. papers

2,481

ext. citations

9

avg. IF

5.14

L-index

#	Paper	IF	Citations
56	The CamSol method of rational design of protein mutants with enhanced solubility. <i>Journal of Molecular Biology</i> , 2015 , 427, 478-90	6.5	208
55	A natural product inhibits the initiation of β -synuclein aggregation and suppresses its toxicity. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017 , 114, E1009-E1017	11.5	177
54	MobiDB 3.0: more annotations for intrinsic disorder, conformational diversity and interactions in proteins. <i>Nucleic Acids Research</i> , 2018 , 46, D471-D476	20.1	143
53	Simultaneous quantification of protein order and disorder. <i>Nature Chemical Biology</i> , 2017 , 13, 339-342	11.7	83
52	Selective targeting of primary and secondary nucleation pathways in A β 2 aggregation using a rational antibody scanning method. <i>Science Advances</i> , 2017 , 3, e1700488	14.3	81
51	Rational design of antibodies targeting specific epitopes within intrinsically disordered proteins. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015 , 112, 9902-7	11.5	78
50	Different soluble aggregates of A β 2 can give rise to cellular toxicity through different mechanisms. <i>Nature Communications</i> , 2019 , 10, 1541	17.4	71
49	A protein homeostasis signature in healthy brains recapitulates tissue vulnerability to Alzheimer's disease. <i>Science Advances</i> , 2016 , 2, e1600947	14.3	68
48	Third generation antibody discovery methods: in silico rational design. <i>Chemical Society Reviews</i> , 2018 , 47, 9137-9157	58.5	64
47	Rapid and accurate in silico solubility screening of a monoclonal antibody library. <i>Scientific Reports</i> , 2017 , 7, 8200	4.9	63
46	Parapred: antibody paratope prediction using convolutional and recurrent neural networks. <i>Bioinformatics</i> , 2018 , 34, 2944-2950	7.2	62
45	The s2D method: simultaneous sequence-based prediction of the statistical populations of ordered and disordered regions in proteins. <i>Journal of Molecular Biology</i> , 2015 , 427, 982-996	6.5	60
44	Targeting disordered proteins with small molecules using entropy. <i>Trends in Biochemical Sciences</i> , 2015 , 40, 491-6	10.3	58
43	Proteome-wide observation of the phenomenon of life on the edge of solubility. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2020 , 117, 1015-1020	11.5	52
42	In vitro and in silico assessment of the developability of a designed monoclonal antibody library. <i>MAbs</i> , 2019 , 11, 388-400	6.6	43
41	Rational design of mutations that change the aggregation rate of a protein while maintaining its native structure and stability. <i>Scientific Reports</i> , 2016 , 6, 25559	4.9	41
40	Understanding the frustration arising from the competition between function, misfolding, and aggregation in a globular protein. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2014 , 111, 14141-6	11.5	36

39	Massively parallel <i>C. elegans</i> tracking provides multi-dimensional fingerprints for phenotypic discovery. <i>Journal of Neuroscience Methods</i> , 2018 , 306, 57-67	3	35
38	Conformational dynamics in crystals reveal the molecular bases for D76N beta-2 microglobulin aggregation propensity. <i>Nature Communications</i> , 2018 , 9, 1658	17.4	35
37	Developability Assessment of Engineered Monoclonal Antibody Variants with a Complex Self-Association Behavior Using Complementary Analytical and in Silico Tools. <i>Molecular Pharmaceutics</i> , 2018 , 15, 5697-5710	5.6	33
36	Oxetane Grafts Installed Site-Selectively on Native Disulfides to Enhance Protein Stability and Activity In Vivo. <i>Angewandte Chemie - International Edition</i> , 2017 , 56, 14963-14967	16.4	32
35	A chemical kinetic basis for measuring translation initiation and elongation rates from ribosome profiling data. <i>PLoS Computational Biology</i> , 2019 , 15, e1007070	5	26
34	Rational design of a conformation-specific antibody for the quantification of A β oligomers. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2020 , 117, 13509-13518	11.5	26
33	Attentive Cross-Modal Paratope Prediction. <i>Journal of Computational Biology</i> , 2019 , 26, 536-545	1.7	24
32	Protein Solubility Predictions Using the CamSol Method in the Study of Protein Homeostasis. <i>Cold Spring Harbor Perspectives in Biology</i> , 2019 , 11,	10.2	22
31	A Rational Design Strategy for the Selective Activity Enhancement of a Molecular Chaperone toward a Target Substrate. <i>Biochemistry</i> , 2015 , 54, 5103-12	3.2	22
30	Supersaturated proteins are enriched at synapses and underlie cell and tissue vulnerability in Alzheimer's disease. <i>Heliyon</i> , 2019 , 5, e02589	3.6	17
29	Fibrillogenic propensity of the GroEL apical domain: a Janus-faced minichaperone. <i>FEBS Letters</i> , 2012 , 586, 1120-7	3.8	13
28	Corneal Dystrophy Mutations Drive Pathogenesis by Targeting TGFBIp Stability and Solubility in a Latent Amyloid-forming Domain. <i>Journal of Molecular Biology</i> , 2018 , 430, 1116-1140	6.5	12
27	Identifying A- and P-site locations on ribosome-protected mRNA fragments using Integer Programming. <i>Scientific Reports</i> , 2019 , 9, 6256	4.9	11
26	A Rationally Designed Hsp70 Variant Rescues the Aggregation-Associated Toxicity of Human IAPP in Cultured Pancreatic Islet β Cells. <i>International Journal of Molecular Sciences</i> , 2018 , 19,	6.3	11
25	Delivery of Native Proteins into <i>C. elegans</i> Using a Transduction Protocol Based on Lipid Vesicles. <i>Scientific Reports</i> , 2017 , 7, 15045	4.9	11
24	MonteGrappa: An iterative Monte Carlo program to optimize biomolecular potentials in simplified models. <i>Computer Physics Communications</i> , 2015 , 186, 93-104	4.2	10
23	Inherent Biophysical Properties Modulate the Toxicity of Soluble Amyloidogenic Light Chains. <i>Journal of Molecular Biology</i> , 2020 , 432, 845-860	6.5	10
22	Quantifying misfolded protein oligomers as drug targets and biomarkers in Alzheimer and Parkinson diseases. <i>Nature Reviews Chemistry</i> , 2021 , 5, 277-294	34.6	10

21	Oxetane Grafts Installed Site-Selectively on Native Disulfides to Enhance Protein Stability and Activity In Vivo. <i>Angewandte Chemie</i> , 2017 , 129, 15159-15163	3.6	8
20	Subdomain architecture and stability of a giant repeat protein. <i>Journal of Physical Chemistry B</i> , 2013 , 117, 13029-37	3.4	8
19	Assessment of Therapeutic Antibody Developability by Combinations of In Vitro and In Silico Methods. <i>Methods in Molecular Biology</i> , 2022 , 2313, 57-113	1.4	8
18	Rationally Designed Antibodies as Research Tools to Study the Structure-Toxicity Relationship of Amyloid- β Oligomers. <i>International Journal of Molecular Sciences</i> , 2020 , 21,	6.3	7
17	Current advances in biopharmaceutical informatics: guidelines, impact and challenges in the computational developability assessment of antibody therapeutics.. <i>MAbs</i> , 2022 , 14, 2020082	6.6	6
16	Biochemical and biophysical comparison of human and mouse beta-2 microglobulin reveals the molecular determinants of low amyloid propensity. <i>FEBS Journal</i> , 2020 , 287, 546-560	5.7	6
15	Systematic Activity Maturation of a Single-Domain Antibody with Non-canonical Amino Acids through Chemical Mutagenesis. <i>Cell Chemical Biology</i> , 2021 , 28, 70-77.e5	8.2	6
14	Comparative Studies in the A30P and A53T β Synuclein Strains to Investigate the Molecular Origins of Parkinson's Disease. <i>Frontiers in Cell and Developmental Biology</i> , 2021 , 9, 552549	5.7	5
13	A method of predicting the in vitro fibril formation propensity of A β 0 mutants based on their inclusion body levels in E. coli. <i>Scientific Reports</i> , 2019 , 9, 3680	4.9	4
12	Iterative derivation of effective potentials to sample the conformational space of proteins at atomistic scale. <i>Journal of Chemical Physics</i> , 2014 , 140, 195101	3.9	4
11	A rationally designed bicyclic peptide remodels A β 2 aggregation in vitro and reduces its toxicity in a worm model of Alzheimer's disease. <i>Scientific Reports</i> , 2020 , 10, 15280	4.9	4
10	Neuroserpin and transthyretin are extracellular chaperones that preferentially inhibit amyloid formation. <i>Science Advances</i> , 2021 , 7, eabf7606	14.3	4
9	Pairs of amino acids at the P- and A-sites of the ribosome predictably and causally modulate translation-elongation rates. <i>Journal of Molecular Biology</i> , 2020 , 432, 166696	6.5	3
8	An open-source automated PEG precipitation assay to measure the relative solubility of proteins with low material requirement. <i>Scientific Reports</i> , 2021 , 11, 21932	4.9	3
7	Fragment-based computational design of antibodies targeting structured epitopes		3
6	Modulating the cardiotoxic behaviour of immunoglobulin light chain dimers through point mutations. <i>Amyloid: the International Journal of Experimental and Clinical Investigation: the Official Journal of the International Society of Amyloidosis</i> , 2019 , 26, 105-106	2.7	2
5	Computational maturation of a single-domain antibody against A β 2 aggregation.. <i>Chemical Science</i> , 2021 , 12, 13940-13948	9.4	2
4	Conformational Entropy as a Potential Liability of Computationally Designed Antibodies. <i>Biomolecules</i> , 2022 , 12, 718	5.9	2

3	Paratope Prediction using Convolutional and Recurrent Neural Networks		1
2	Rationally Designed Bicyclic Peptides Prevent the Conversion of A β 2 Assemblies Into Fibrillar Structures. <i>Frontiers in Neuroscience</i> , 2021 , 15, 623097	5.1	1
1	An aggregation inhibitor specific to oligomeric intermediates of A β 2 derived from phage display libraries of stable, small proteins.. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2022 , 119, e2121966119	11.5	1